



Full wwPDB NMR Structure Validation Report ⓘ

Jul 18, 2023 – 02:06 PM EDT

PDB ID : 8DSX
BMRB ID : 51520
Title : NMR STRUCTURE OF THE BACTERIOPHAGE LAMBDA EA22 C-
TERMINAL DOMAIN
Authors : Donaldson, L.W.
Deposited on : 2022-07-23

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
wwPDB-RCI : v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV : Wang et al. (2010)
wwPDB-ShiftChecker : v1.2
BMRB Restraints Analysis : v1.2
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.34

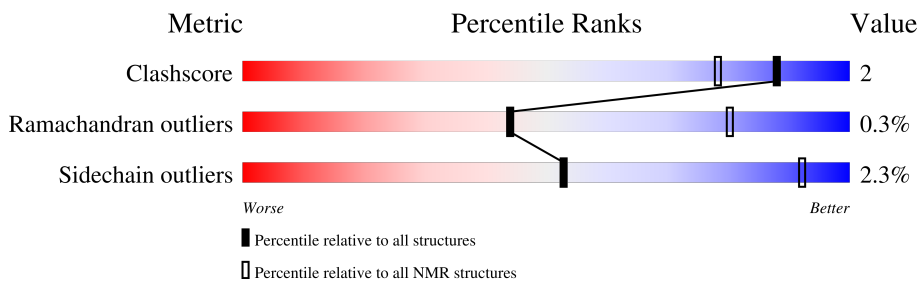
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment is 41%.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	158937	12864
Ramachandran outliers	154571	11451
Sidechain outliers	154315	11428

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	93	
1	B	93	

2 Ensemble composition and analysis i

This entry contains 20 models. Model 5 is the overall representative, medoid model (most similar to other models). The authors have identified model 1 as representative, based on the following criterion: *lowest energy*.

The following residues are included in the computation of the global validation metrics.

Well-defined (core) protein residues			
Well-defined core	Residue range (total)	Backbone RMSD (Å)	Medoid model
1	A:18-A:58, A:65-A:77, B:118-B:158, B:165-B:177 (108)	0.95	5

Ill-defined regions of proteins are excluded from the global statistics.

Ligands and non-protein polymers are included in the analysis.

The models can be grouped into 4 clusters and 3 single-model clusters were found.

Cluster number	Models
1	2, 4, 5, 6, 9, 10, 11, 13, 17, 20
2	1, 7, 12
3	16, 18
4	3, 15
Single-model clusters	8; 14; 19

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2081 atoms, of which 1025 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called Protein ea22.

Mol	Chain	Residues	Atoms					Trace	
			Total	C	H	N	O		S
1	A	65	1040	337	512	91	97	3	0
1	B	65	1041	337	513	91	97	3	0

There are 24 discrepancies between the modelled and reference sequences:

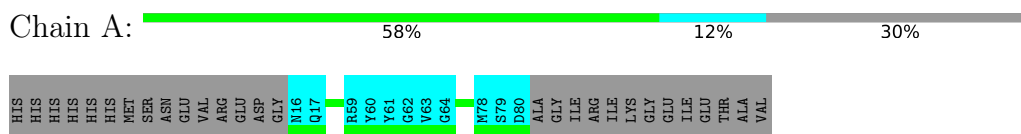
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	HIS	-	expression tag	UNP P03756
A	2	HIS	-	expression tag	UNP P03756
A	3	HIS	-	expression tag	UNP P03756
A	4	HIS	-	expression tag	UNP P03756
A	5	HIS	-	expression tag	UNP P03756
A	6	HIS	-	expression tag	UNP P03756
A	7	MET	-	expression tag	UNP P03756
A	89	ILE	-	expression tag	UNP P03756
A	90	GLU	-	expression tag	UNP P03756
A	91	THR	-	expression tag	UNP P03756
A	92	ALA	-	expression tag	UNP P03756
A	93	VAL	-	expression tag	UNP P03756
B	101	HIS	-	expression tag	UNP P03756
B	102	HIS	-	expression tag	UNP P03756
B	103	HIS	-	expression tag	UNP P03756
B	104	HIS	-	expression tag	UNP P03756
B	105	HIS	-	expression tag	UNP P03756
B	106	HIS	-	expression tag	UNP P03756
B	107	MET	-	expression tag	UNP P03756
B	189	ILE	-	expression tag	UNP P03756
B	190	GLU	-	expression tag	UNP P03756
B	191	THR	-	expression tag	UNP P03756
B	192	ALA	-	expression tag	UNP P03756
B	193	VAL	-	expression tag	UNP P03756

4 Residue-property plots [i](#)

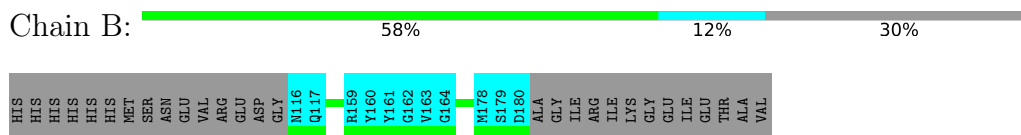
4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: Protein ea22



- Molecule 1: Protein ea22

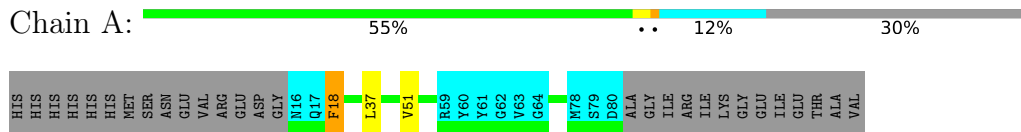


4.2 Scores per residue for each member of the ensemble

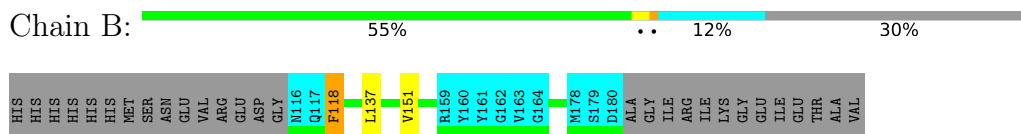
Colouring as in section 4.1 above.

4.2.1 Score per residue for model 1

- Molecule 1: Protein ea22

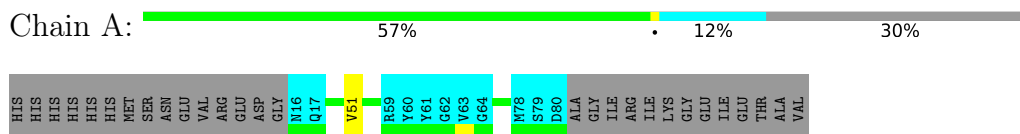


- Molecule 1: Protein ea22

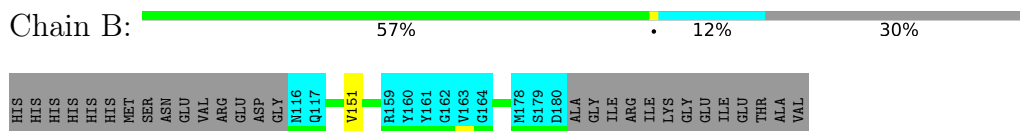


4.2.2 Score per residue for model 2

- Molecule 1: Protein ea22

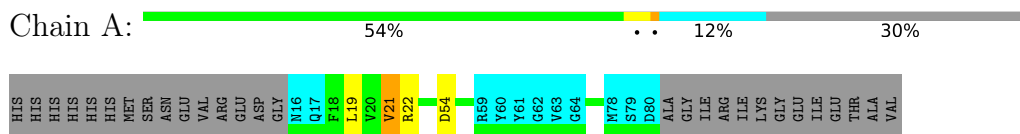


- Molecule 1: Protein ea22

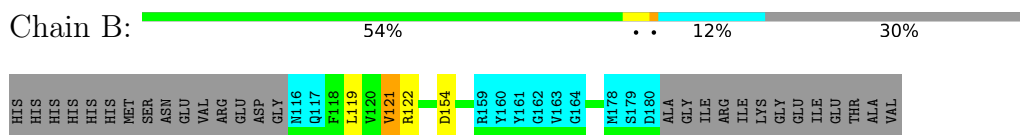


4.2.3 Score per residue for model 3

- Molecule 1: Protein ea22

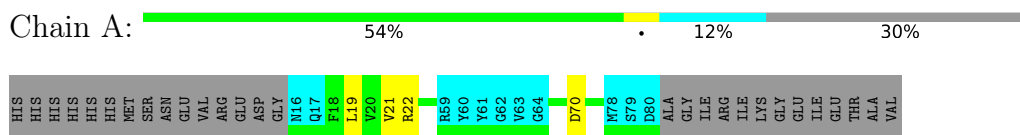


- Molecule 1: Protein ea22

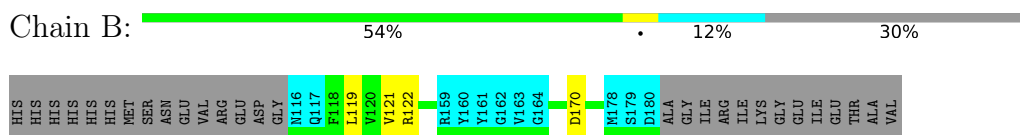


4.2.4 Score per residue for model 4

- Molecule 1: Protein ea22

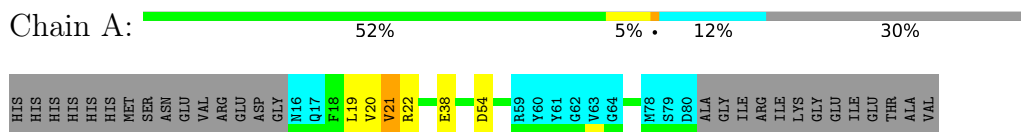


- Molecule 1: Protein ea22

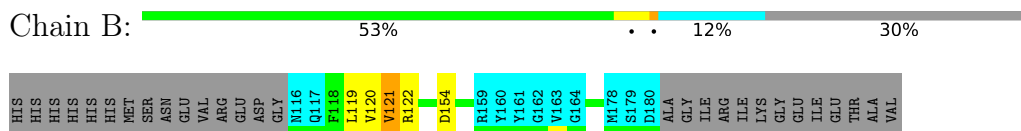


4.2.5 Score per residue for model 5 (medoid)

- Molecule 1: Protein ea22

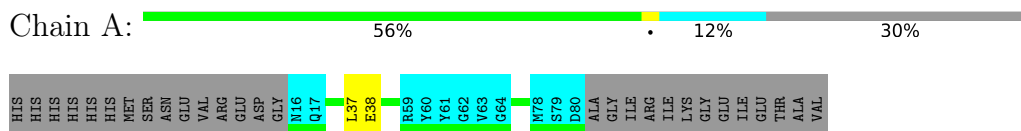


- Molecule 1: Protein ea22

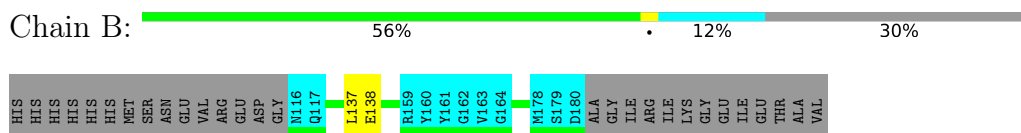


4.2.6 Score per residue for model 6

- Molecule 1: Protein ea22

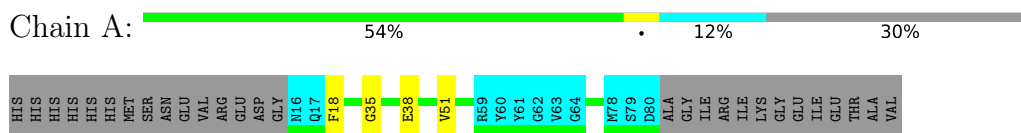


- Molecule 1: Protein ea22

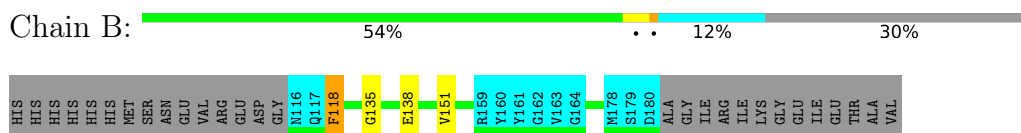


4.2.7 Score per residue for model 7

- Molecule 1: Protein ea22

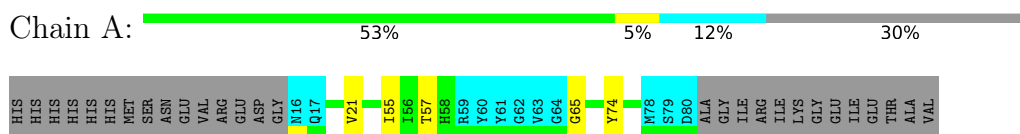


- Molecule 1: Protein ea22

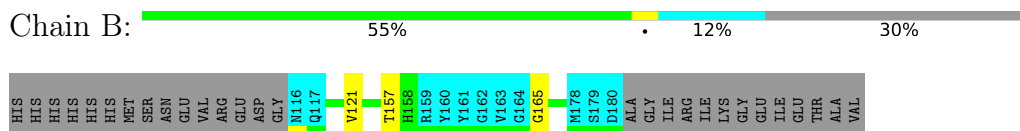


4.2.8 Score per residue for model 8

- Molecule 1: Protein ea22

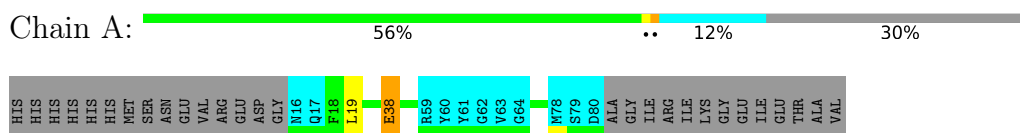


- Molecule 1: Protein ea22

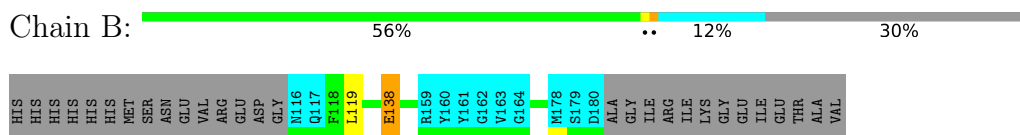


4.2.9 Score per residue for model 9

- Molecule 1: Protein ea22

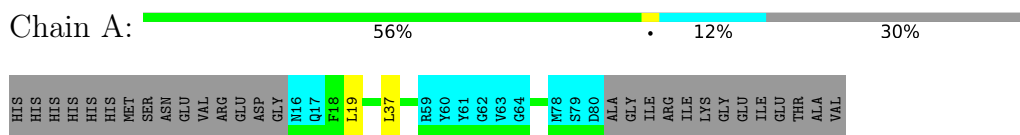


- Molecule 1: Protein ea22



4.2.10 Score per residue for model 10

- Molecule 1: Protein ea22

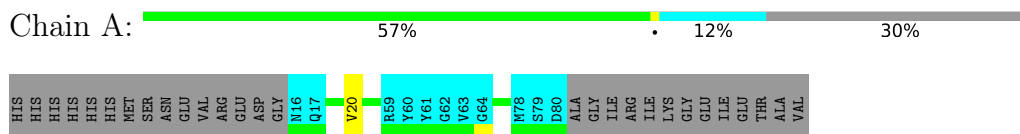


- Molecule 1: Protein ea22

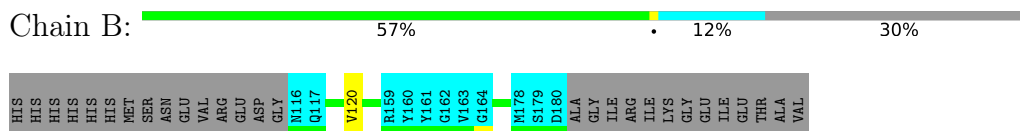


4.2.11 Score per residue for model 11

- Molecule 1: Protein ea22

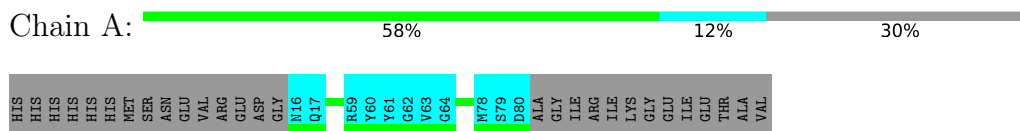


- Molecule 1: Protein ea22

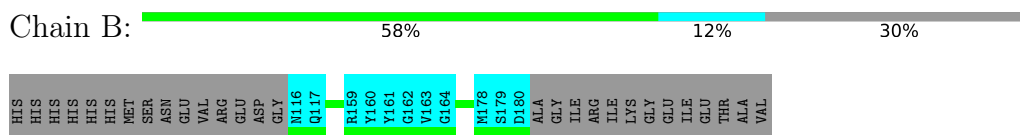


4.2.12 Score per residue for model 12

- Molecule 1: Protein ea22

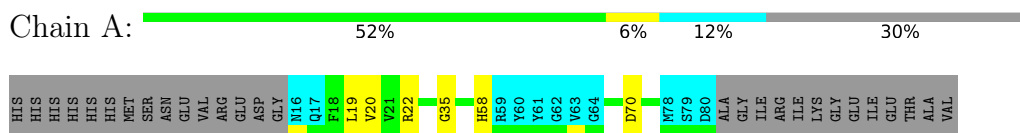


- Molecule 1: Protein ea22

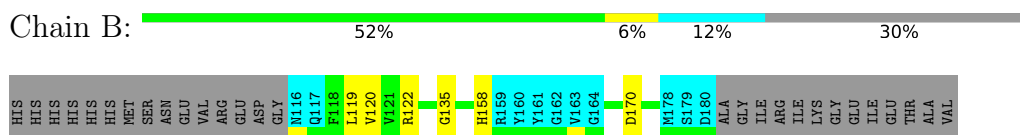


4.2.13 Score per residue for model 13

- Molecule 1: Protein ea22



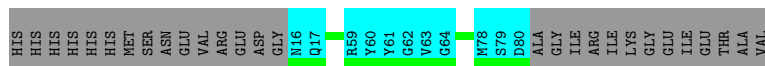
- Molecule 1: Protein ea22



4.2.14 Score per residue for model 14

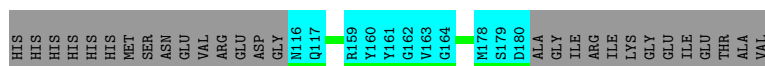
- Molecule 1: Protein ea22

Chain A:  58% 12% 30%



- Molecule 1: Protein ea22

Chain B:  58% 12% 30%



4.2.15 Score per residue for model 15

- Molecule 1: Protein ea22

Chain A:  55% 12% 30%



- Molecule 1: Protein ea22

Chain B:  55% 12% 30%



4.2.16 Score per residue for model 16

- Molecule 1: Protein ea22

Chain A:  56% 12% 30%



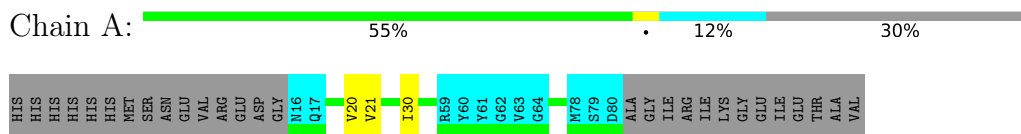
- Molecule 1: Protein ea22

Chain B:  56% 12% 30%

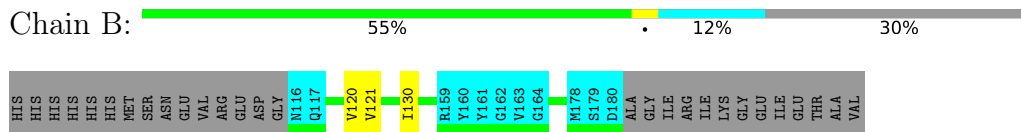


4.2.17 Score per residue for model 17

- Molecule 1: Protein ea22

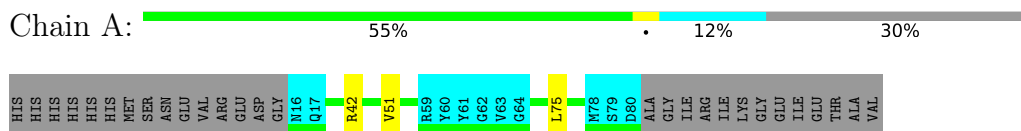


- Molecule 1: Protein ea22

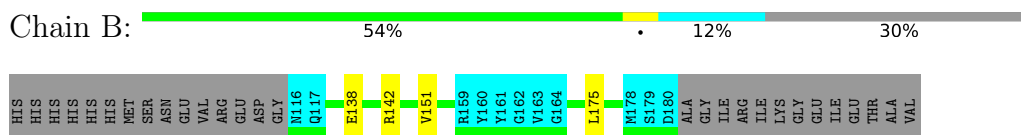


4.2.18 Score per residue for model 18

- Molecule 1: Protein ea22

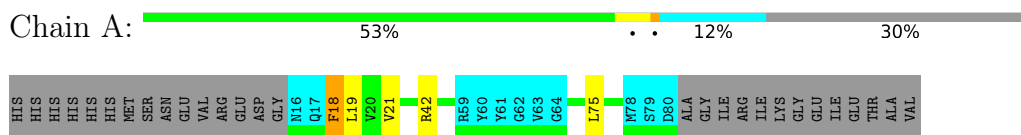


- Molecule 1: Protein ea22

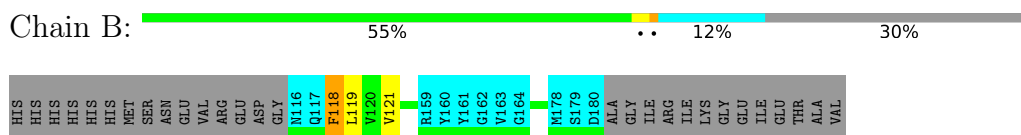


4.2.19 Score per residue for model 19

- Molecule 1: Protein ea22

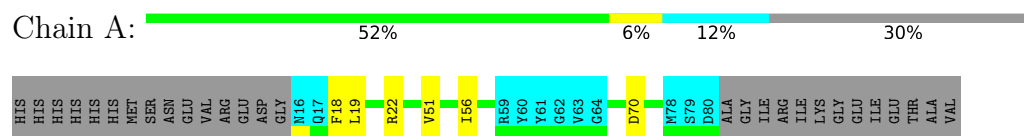


- Molecule 1: Protein ea22

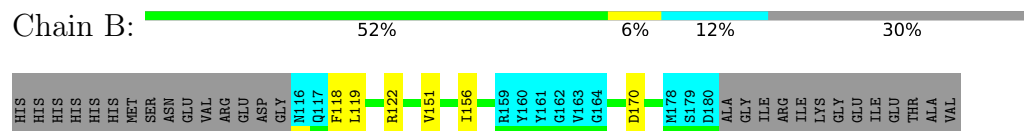


4.2.20 Score per residue for model 20

- Molecule 1: Protein ea22



- Molecule 1: Protein ea22



5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing, simulated annealing*.

Of the 1000 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CYANA	refinement	3
Rosetta	refinement	2021.16.61629
CYANA	structure calculation	3

The following table shows chemical shift validation statistics as aggregates over all chemical shift files. Detailed validation can be found in section 7 of this report.

Chemical shift file(s)	working_cs.cif
Number of chemical shift lists	1
Total number of shifts	798
Number of shifts mapped to atoms	685
Number of unparsed shifts	0
Number of shifts with mapping errors	113
Number of shifts with mapping warnings	0
Assignment completeness (well-defined parts)	41%

6 Model quality i

6.1 Standard geometry i

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts i

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	439	435	435	2±1
1	B	439	435	435	2±1
All	All	17560	17400	17400	67

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

All unique clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:A:48:ASP:O	1:A:51:VAL:HG22	0.71	1.86	15	1
1:B:148:ASP:O	1:B:151:VAL:HG22	0.69	1.86	15	1
1:A:54:ASP:OD1	1:B:122:ARG:NH1	0.63	2.31	3	2
1:A:22:ARG:NH1	1:B:154:ASP:OD1	0.62	2.31	3	2
1:B:137:LEU:HD13	1:B:137:LEU:C	0.62	2.15	10	3
1:A:37:LEU:HD13	1:A:37:LEU:C	0.61	2.15	16	3
1:A:22:ARG:NH2	1:B:170:ASP:OD2	0.58	2.36	20	3
1:B:137:LEU:HD13	1:B:137:LEU:O	0.58	1.98	16	2
1:A:37:LEU:HD13	1:A:37:LEU:O	0.57	1.98	16	3
1:A:70:ASP:OD2	1:B:122:ARG:NH2	0.57	2.36	20	3
1:B:138:GLU:OE1	1:B:138:GLU:N	0.56	2.36	7	4
1:B:118:PHE:C	1:B:118:PHE:CD1	0.56	2.78	1	2
1:A:38:GLU:OE1	1:A:38:GLU:N	0.56	2.36	7	4

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Atom-1	Atom-2	Clash(Å)	Distance(Å)	Models	
				Worst	Total
1:B:118:PHE:C	1:B:118:PHE:CD2	0.55	2.80	20	2
1:A:20:VAL:HG22	1:A:30:ILE:CD1	0.55	2.32	17	1
1:B:120:VAL:HG22	1:B:130:ILE:CD1	0.54	2.32	17	1
1:A:18:PHE:C	1:A:18:PHE:CD2	0.54	2.80	20	1
1:A:18:PHE:C	1:A:18:PHE:CD1	0.54	2.78	1	2
1:B:137:LEU:O	1:B:137:LEU:HD13	0.52	2.04	1	1
1:A:21:VAL:HG22	1:A:21:VAL:O	0.52	2.05	3	2
1:B:121:VAL:HG22	1:B:121:VAL:O	0.51	2.04	5	2
1:A:37:LEU:HD23	1:A:37:LEU:O	0.48	2.09	6	1
1:B:137:LEU:O	1:B:137:LEU:HD23	0.47	2.09	6	1
1:B:137:LEU:C	1:B:137:LEU:CD1	0.47	2.83	16	2
1:A:37:LEU:C	1:A:37:LEU:CD1	0.46	2.83	16	2
1:A:38:GLU:CD	1:A:38:GLU:H	0.45	2.14	9	1
1:A:20:VAL:HG22	1:A:30:ILE:HD12	0.45	1.87	17	1
1:B:120:VAL:HG22	1:B:130:ILE:HD12	0.45	1.87	17	1
1:B:138:GLU:H	1:B:138:GLU:CD	0.45	2.14	9	1
1:A:42:ARG:CG	1:A:75:LEU:HD21	0.42	2.45	18	2
1:B:142:ARG:CG	1:B:175:LEU:HD21	0.42	2.45	18	1
1:B:123:HIS:HA	1:B:151:VAL:HG12	0.42	1.91	15	1
1:B:120:VAL:O	1:B:120:VAL:HG13	0.42	2.15	5	2
1:A:20:VAL:O	1:A:20:VAL:HG13	0.41	2.15	13	3
1:A:23:HIS:HA	1:A:51:VAL:HG12	0.41	1.91	15	1
1:B:120:VAL:HG13	1:B:120:VAL:O	0.41	2.16	11	1
1:A:55:ILE:CD1	1:A:74:TYR:CD2	0.40	3.05	8	1

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	54/93 (58%)	52±1 (97±2%)	1±1 (3±2%)	0±0 (0±1%)	44	80
1	B	54/93 (58%)	52±1 (97±2%)	1±1 (3±2%)	0±0 (0±1%)	44	80
All	All	2160/3720 (58%)	2096 (97%)	58 (3%)	6 (0%)	44	80

All 4 unique Ramachandran outliers are listed below. They are sorted by the frequency of occur-

rence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	35	GLY	2
1	B	135	GLY	2
1	A	65	GLY	1
1	B	165	GLY	1

6.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	49/81 (60%)	48±1 (98±2%)	1±1 (2±2%)	53	92
1	B	49/81 (60%)	48±1 (98±2%)	1±1 (2±2%)	53	92
All	All	1960/3240 (60%)	1914 (98%)	46 (2%)	53	92

All 10 unique residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type	Models (Total)
1	A	19	LEU	8
1	B	119	LEU	8
1	A	21	VAL	6
1	B	121	VAL	6
1	A	51	VAL	5
1	B	151	VAL	5
1	A	18	PHE	3
1	B	118	PHE	3
1	A	38	GLU	1
1	B	138	GLU	1

6.3.3 RNA [i](#)

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation i

The completeness of assignment taking into account all chemical shift lists is 41% for the well-defined parts and 37% for the entire structure.

7.1 Chemical shift list 1

File name: working_cs.cif

Chemical shift list name: *assigned_chemical_shifts_1*

7.1.1 Bookkeeping i

The following table shows the results of parsing the chemical shift list and reports the number of nuclei with statistically unusual chemical shifts.

Total number of shifts	798
Number of shifts mapped to atoms	685
Number of unparsed shifts	0
Number of shifts with mapping errors	113
Number of shifts with mapping warnings	0
Number of shift outliers (ShiftChecker)	0

The following errors were found when reading this chemical shift list.

- Chemical shift has been reported more than once. All 4 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	93	VAL	HG12	0.879	0.000	2
1	A	93	VAL	HG13	0.879	0.000	2
1	A	93	VAL	HG22	0.905	0.000	2
1	A	93	VAL	HG23	0.905	0.000	2

The following assigned chemical shifts were not mapped to the molecules present in the coordinate file.

- No matching atom found in the structure. All 113 occurrences are reported below.

List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	81	ALA	H	8.04	0.004	1
1	A	81	ALA	HA	4.21	0.005	1
1	A	81	ALA	HB1	1.409	0.006	1
1	A	81	ALA	HB2	1.409	0.006	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	81	ALA	HB3	1.409	0.006	1
1	A	81	ALA	C	178.631	0.000	1
1	A	81	ALA	CA	53.271	0.077	1
1	A	81	ALA	CB	18.821	0.136	1
1	A	81	ALA	N	123.401	0.000	1
1	A	82	GLY	H	8.193	0.003	1
1	A	82	GLY	C	174.525	0.014	1
1	A	82	GLY	CA	45.494	0.029	1
1	A	82	GLY	N	106.481	0.000	1
1	A	83	ILE	H	7.741	0.004	1
1	A	83	ILE	HG21	0.897	0.000	1
1	A	83	ILE	HG22	0.897	0.000	1
1	A	83	ILE	HG23	0.897	0.000	1
1	A	83	ILE	HD11	0.858	0.000	1
1	A	83	ILE	HD12	0.858	0.000	1
1	A	83	ILE	HD13	0.858	0.000	1
1	A	83	ILE	C	176.402	0.056	1
1	A	83	ILE	CA	61.382	0.000	1
1	A	83	ILE	CB	38.453	0.000	1
1	A	83	ILE	CG2	17.51	0.000	1
1	A	83	ILE	CD1	13.017	0.000	1
1	A	83	ILE	N	119.479	0.000	1
1	A	84	ARG	H	8.148	0.001	1
1	A	84	ARG	C	176.331	0.042	1
1	A	84	ARG	CA	56.318	0.000	1
1	A	84	ARG	CB	30.619	0.000	1
1	A	84	ARG	N	123.917	0.000	1
1	A	85	ILE	H	8.059	0.005	1
1	A	85	ILE	HG21	0.897	0.000	1
1	A	85	ILE	HG22	0.897	0.000	1
1	A	85	ILE	HG23	0.897	0.000	1
1	A	85	ILE	HD11	0.854	0.000	1
1	A	85	ILE	HD12	0.854	0.000	1
1	A	85	ILE	HD13	0.854	0.000	1
1	A	85	ILE	C	176.317	0.000	1
1	A	85	ILE	CA	61.143	0.048	1
1	A	85	ILE	CB	38.453	0.000	1
1	A	85	ILE	CG2	17.466	0.000	1
1	A	85	ILE	CD1	12.935	0.000	1
1	A	85	ILE	N	121.73	0.000	1
1	A	86	LYS	H	8.277	0.005	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	86	LYS	C	176.811	0.042	1
1	A	86	LYS	CA	56.617	0.084	1
1	A	86	LYS	CB	32.924	0.144	1
1	A	86	LYS	CG	24.933	0.000	1
1	A	86	LYS	CD	29.412	0.000	1
1	A	86	LYS	N	125.108	0.000	1
1	A	87	GLY	H	8.269	0.004	1
1	A	87	GLY	C	173.878	0.012	1
1	A	87	GLY	CA	45.156	0.046	1
1	A	87	GLY	N	109.835	0.014	1
1	A	88	GLU	H	8.201	0.003	1
1	A	88	GLU	C	176.486	0.000	1
1	A	88	GLU	CA	56.546	0.000	1
1	A	88	GLU	CB	30.421	0.000	1
1	A	88	GLU	CG	36.286	0.000	1
1	A	88	GLU	N	120.507	0.000	1
1	A	89	ILE	H	8.109	0.003	1
1	A	89	ILE	HG21	0.902	0.000	1
1	A	89	ILE	HG22	0.902	0.000	1
1	A	89	ILE	HG23	0.902	0.000	1
1	A	89	ILE	HD11	0.853	0.000	1
1	A	89	ILE	HD12	0.853	0.000	1
1	A	89	ILE	HD13	0.853	0.000	1
1	A	89	ILE	C	176.402	0.000	1
1	A	89	ILE	CG2	17.466	0.000	1
1	A	89	ILE	CD1	12.864	0.000	1
1	A	89	ILE	N	120.957	0.000	1
1	A	90	GLU	H	8.463	0.001	1
1	A	90	GLU	C	176.345	0.000	1
1	A	90	GLU	CA	56.623	0.082	1
1	A	90	GLU	CB	30.329	0.092	1
1	A	90	GLU	CG	36.494	0.000	1
1	A	90	GLU	N	124.985	0.000	1
1	A	91	THR	H	8.098	0.004	1
1	A	91	THR	HA	4.321	0.000	1
1	A	91	THR	HB	4.184	0.001	1
1	A	91	THR	HG21	1.203	0.000	1
1	A	91	THR	HG22	1.203	0.000	1
1	A	91	THR	HG23	1.203	0.000	1
1	A	91	THR	C	173.904	0.014	1
1	A	91	THR	CA	61.744	0.126	1

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List ID	Chain	Res	Type	Atom	Shift Data		
					Value	Uncertainty	Ambiguity
1	A	91	THR	CB	69.959	0.047	1
1	A	91	THR	CG2	21.683	0.096	1
1	A	91	THR	N	115.729	0.000	1
1	A	92	ALA	H	8.281	0.003	1
1	A	92	ALA	HA	4.414	0.000	1
1	A	92	ALA	HB1	1.393	0.004	1
1	A	92	ALA	HB2	1.393	0.004	1
1	A	92	ALA	HB3	1.393	0.004	1
1	A	92	ALA	C	176.571	0.000	1
1	A	92	ALA	CA	52.537	0.030	1
1	A	92	ALA	CB	19.289	0.078	1
1	A	92	ALA	N	127.675	0.046	1
1	A	93	VAL	H	7.639	0.003	1
1	A	93	VAL	HA	4.043	0.000	1
1	A	93	VAL	HB	2.091	0.000	1
1	A	93	VAL	HG11	0.879	0.000	2
1	A	93	VAL	HG12	0.879	0.000	2
1	A	93	VAL	HG13	0.879	0.000	2
1	A	93	VAL	HG21	0.905	0.000	2
1	A	93	VAL	HG22	0.905	0.000	2
1	A	93	VAL	HG23	0.905	0.000	2
1	A	93	VAL	C	180.917	0.000	1
1	A	93	VAL	CA	63.566	0.000	1
1	A	93	VAL	CB	32.914	0.000	1
1	A	93	VAL	CG1	20.033	0.000	2
1	A	93	VAL	CG2	21.554	0.000	2
1	A	93	VAL	N	123.346	0.000	1

7.1.2 Chemical shift referencing [i](#)

The following table shows the suggested chemical shift referencing corrections.

Nucleus	# values	Correction \pm precision, ppm	Suggested action
$^{13}\text{C}_\alpha$	66	0.23 ± 0.19	None needed (< 0.5 ppm)
$^{13}\text{C}_\beta$	59	-0.09 ± 0.29	None needed (< 0.5 ppm)
$^{13}\text{C}'$	66	0.07 ± 0.25	None needed (< 0.5 ppm)
^{15}N	59	0.56 ± 0.62	None needed (imprecise)

7.1.3 Completeness of resonance assignments [i](#)

The following table shows the completeness of the chemical shift assignments for the well-defined regions of the structure. The overall completeness is 41%, i.e. 636 atoms were assigned a chemical shift out of a possible 1536. 0 out of 22 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	231/536 (43%)	91/218 (42%)	97/216 (45%)	43/102 (42%)
Sidechain	366/862 (42%)	250/564 (44%)	112/274 (41%)	4/24 (17%)
Aromatic	39/138 (28%)	21/72 (29%)	17/56 (30%)	1/10 (10%)
Overall	636/1536 (41%)	362/854 (42%)	226/546 (41%)	48/136 (35%)

The following table shows the completeness of the chemical shift assignments for the full structure. The overall completeness is 37%, i.e. 681 atoms were assigned a chemical shift out of a possible 1822. 0 out of 24 assigned methyl groups (LEU and VAL) were assigned stereospecifically.

	Total	¹H	¹³C	¹⁵N
Backbone	253/650 (39%)	100/266 (38%)	107/260 (41%)	46/124 (37%)
Sidechain	389/998 (39%)	266/650 (41%)	119/314 (38%)	4/34 (12%)
Aromatic	39/174 (22%)	21/88 (24%)	17/76 (22%)	1/10 (10%)
Overall	681/1822 (37%)	387/1004 (39%)	243/650 (37%)	51/168 (30%)

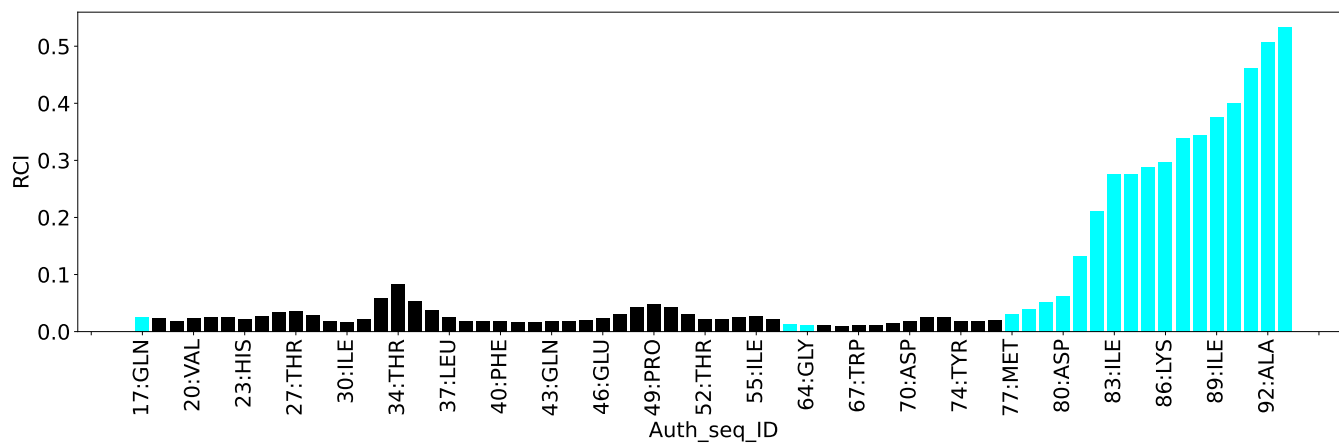
7.1.4 Statistically unusual chemical shifts [i](#)

There are no statistically unusual chemical shifts.

7.1.5 Random Coil Index (RCI) plots [i](#)

The image below reports *random coil index* values for the protein chains in the structure. The height of each bar gives a probability of a given residue to be disordered, as predicted from the available chemical shifts and the amino acid sequence. A value above 0.2 is an indication of significant predicted disorder. The colour of the bar shows whether the residue is in the well-defined core (black) or in the ill-defined residue ranges (cyan), as described in section 2 on ensemble composition. If well-defined core and ill-defined regions are not identified then it is shown as gray bars.

Random coil index (RCI) for chain A:



8 NMR restraints analysis

8.1 Conformationally restricting restraints

The following table provides the summary of experimentally observed NMR restraints in different categories. Restraints are classified into different categories based on the sequence separation of the atoms involved.

Description	Value
Total distance restraints	904
Intra-residue ($ i-j =0$)	407
Sequential ($ i-j =1$)	159
Medium range ($ i-j >1$ and $ i-j <5$)	79
Long range ($ i-j \geq 5$)	199
Inter-chain	24
Hydrogen bond restraints	36
Disulfide bond restraints	0
Total dihedral-angle restraints	0
Number of unmapped restraints	0
Number of restraints per residue	4.9
Number of long range restraints per residue ¹	1.2

¹Long range hydrogen bonds and disulfide bonds are counted as long range restraints while calculating the number of long range restraints per residue

8.2 Residual restraint violations

This section provides the overview of the restraint violations analysis. The violations are binned as small, medium and large violations based on its absolute value. Average number of violations per model is calculated by dividing the total number of violations in each bin by the size of the ensemble.

8.2.1 Average number of distance violations per model

Distance violations less than 0.1 Å are not included in the calculation.

Bins (Å)	Average number of violations per model	Max (Å)
0.1-0.2 (Small)	4.9	0.2
0.2-0.5 (Medium)	1.4	0.49
>0.5 (Large)	0.2	1.19

8.2.2 Average number of dihedral-angle violations per model [i](#)

Dihedral-angle violations less than 1° are not included in the calculation. There are no dihedral-angle violations

9 Distance violation analysis [i](#)

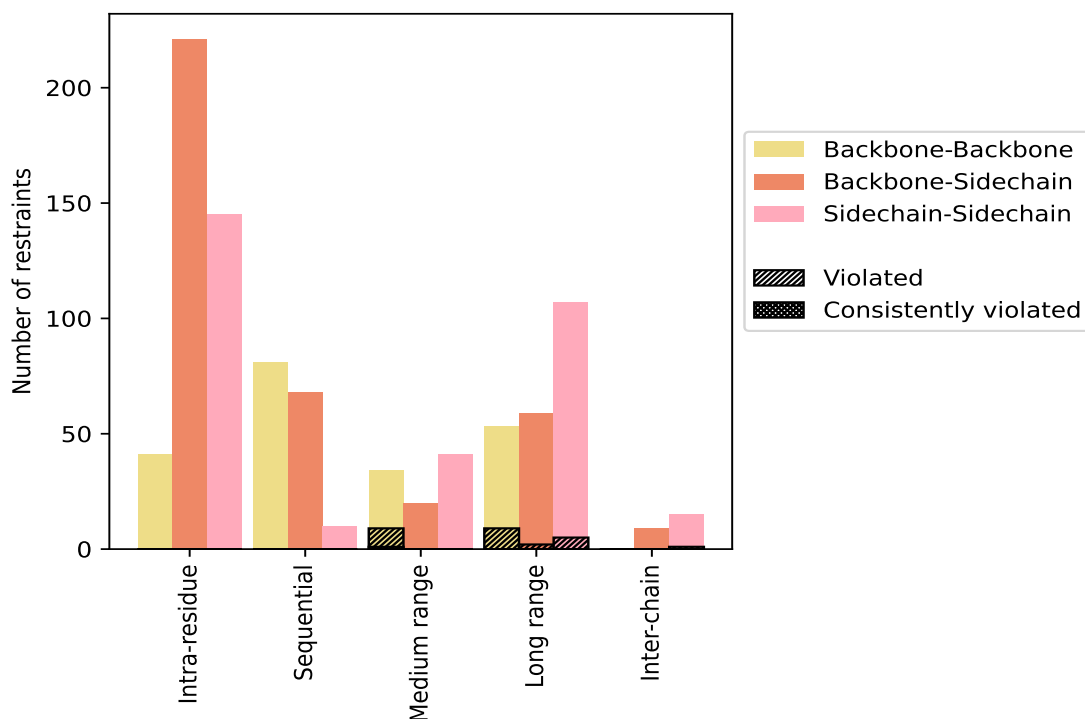
9.1 Summary of distance violations [i](#)

The following table shows the summary of distance violations in different restraint categories based on the sequence separation of the atoms involved. Each category is further sub-divided into three sub-categories based on the atoms involved. Violations less than 0.1 Å are not included in the statistics.

Restrains type	Count	% ¹	Violated ³			Consistently Violated ⁴		
			Count	% ²	% ¹	Count	% ²	% ¹
Intra-residue (i-j =0)	407	45.0	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	41	4.5	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	221	24.4	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	145	16.0	0	0.0	0.0	0	0.0	0.0
Sequential (i-j =1)	159	17.6	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	81	9.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	68	7.5	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	10	1.1	0	0.0	0.0	0	0.0	0.0
Medium range (i-j >1 & i-j <5)	79	8.7	0	0.0	0.0	0	0.0	0.0
Backbone-Backbone	18	2.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	20	2.2	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	41	4.5	0	0.0	0.0	0	0.0	0.0
Long range (i-j ≥5)	199	22.0	7	3.5	0.8	0	0.0	0.0
Backbone-Backbone	33	3.7	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	59	6.5	2	3.4	0.2	0	0.0	0.0
Sidechain-Sidechain	107	11.8	5	4.7	0.6	0	0.0	0.0
Inter-chain	24	2.7	1	4.2	0.1	0	0.0	0.0
Backbone-Backbone	0	0.0	0	0.0	0.0	0	0.0	0.0
Backbone-Sidechain	9	1.0	0	0.0	0.0	0	0.0	0.0
Sidechain-Sidechain	15	1.7	1	6.7	0.1	0	0.0	0.0
Hydrogen bond	36	4.0	18	50.0	2.0	1	2.8	0.1
Disulfide bond	0	0.0	0	0.0	0.0	0	0.0	0.0
Total	904	100.0	26	2.9	2.9	1	0.1	0.1
Backbone-Backbone	209	23.1	18	8.6	2.0	1	0.5	0.1
Backbone-Sidechain	377	41.7	2	0.5	0.2	0	0.0	0.0
Sidechain-Sidechain	318	35.2	6	1.9	0.7	0	0.0	0.0

¹ percentage calculated with respect to the total number of distance restraints, ² percentage calculated with respect to the number of restraints in a particular restraint category, ³ violated in at least one model, ⁴ violated in all the models

9.1.1 Bar chart : Distribution of distance restraints and violations [i](#)



Violated and consistently violated restraints are shown using different hatch patterns in their respective categories. The hydrogen bonds and disulfid bonds are counted in their appropriate category on the x-axis

9.2 Distance violation statistics for each model [i](#)

The following table provides the distance violation statistics for each model in the ensemble. Violations less than 0.1 Å are not included in the statistics.

Model ID	Number of violations						Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total				
1	0	0	5	3	0	8	0.18	0.31	0.06	0.16
2	0	0	2	2	0	4	0.24	0.53	0.17	0.16
3	0	0	4	2	1	7	0.24	0.66	0.18	0.15
4	0	0	2	2	0	4	0.14	0.16	0.02	0.13
5	0	0	3	2	0	5	0.26	0.72	0.23	0.18
6	0	0	3	3	1	7	0.32	1.19	0.36	0.16
7	0	0	2	4	0	6	0.19	0.27	0.05	0.18
8	0	0	2	4	0	6	0.18	0.22	0.04	0.2
9	0	0	3	3	0	6	0.14	0.19	0.03	0.12
10	0	0	2	6	0	8	0.14	0.19	0.03	0.14
11	0	0	2	3	0	5	0.2	0.49	0.15	0.12

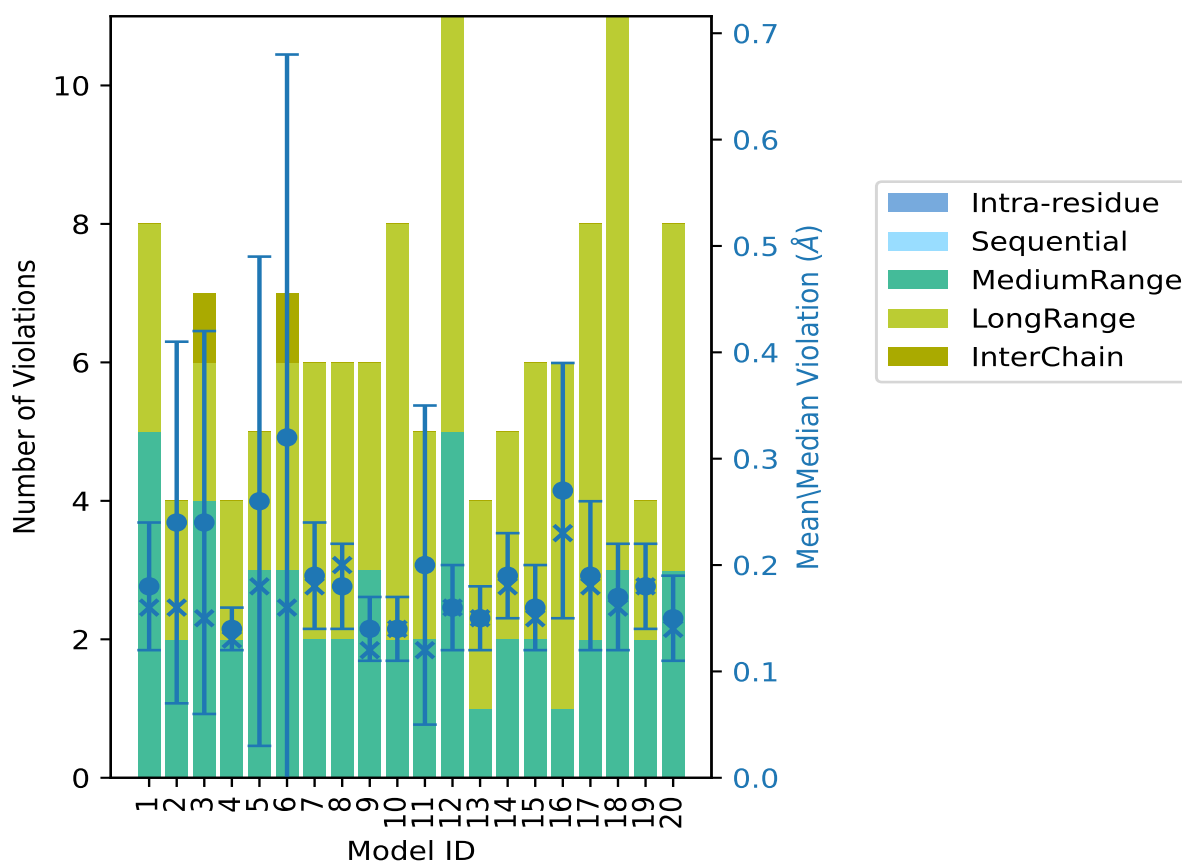
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Model ID	Number of violations					Total	Mean (Å)	Max (Å)	SD ⁶ (Å)	Median (Å)
	IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵					
12	0	0	5	6	0	11	0.16	0.26	0.04	0.16
13	0	0	1	3	0	4	0.15	0.19	0.03	0.15
14	0	0	2	3	0	5	0.19	0.26	0.04	0.18
15	0	0	2	4	0	6	0.16	0.25	0.04	0.15
16	0	0	1	5	0	6	0.27	0.49	0.12	0.23
17	0	0	2	6	0	8	0.19	0.31	0.07	0.18
18	0	0	3	8	0	11	0.17	0.28	0.05	0.16
19	0	0	2	2	0	4	0.18	0.24	0.04	0.18
20	0	0	3	5	0	8	0.15	0.24	0.04	0.14

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints, ⁵Inter-chain restraints, ⁶Standard deviation

9.2.1 Bar graph : Distance Violation statistics for each model [\(i\)](#)



The mean(dot),median(x) and the standard deviation are shown in blue with respect to the y axis on the right

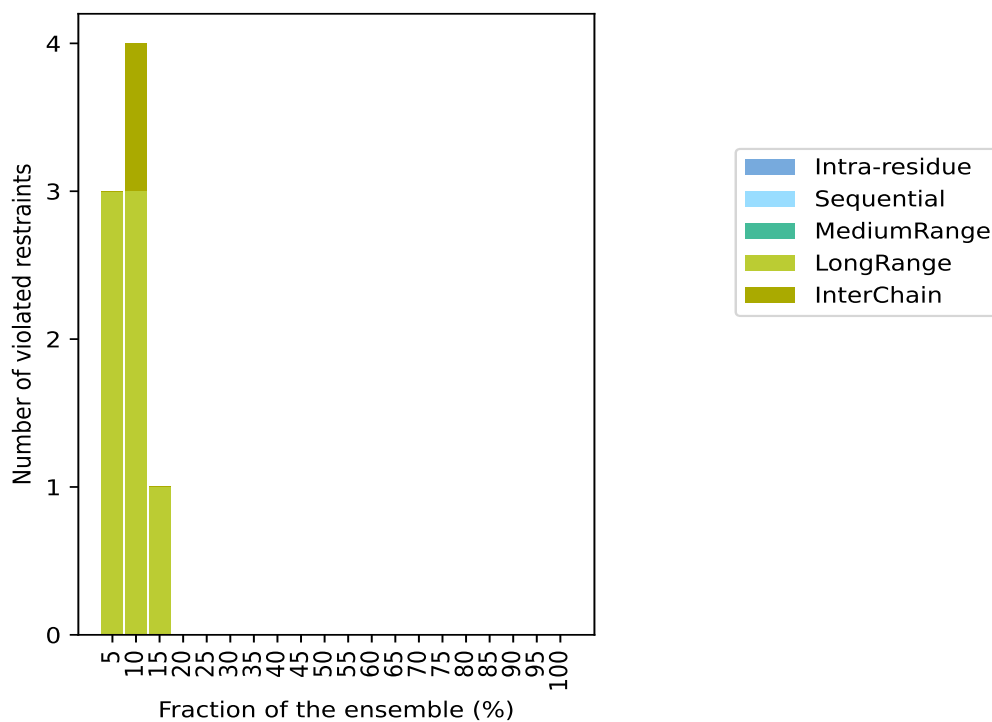
9.3 Distance violation statistics for the ensemble

Violation analysis may find that some restraints are violated in few models and some are violated in most of models. The following table provides this information as number of violated restraints for a given fraction of the ensemble. In total, 860(IR:407, SQ:159, MR:79, LR:192, IC:23) restraints are not violated in the ensemble.

Number of violated restraints						Fraction of the ensemble	
IR ¹	SQ ²	MR ³	LR ⁴	IC ⁵	Total	Count ⁶	%
0	0	0	3	0	3	1	5.0
0	0	0	3	1	4	2	10.0
0	0	0	1	0	1	3	15.0
0	0	0	0	0	0	4	20.0
0	0	0	0	0	0	5	25.0
0	0	0	0	0	0	6	30.0
0	0	0	0	0	0	7	35.0
0	0	0	0	0	0	8	40.0
0	0	0	0	0	0	9	45.0
0	0	0	0	0	0	10	50.0
0	0	0	0	0	0	11	55.0
0	0	0	0	0	0	12	60.0
0	0	0	0	0	0	13	65.0
0	0	0	0	0	0	14	70.0
0	0	0	0	0	0	15	75.0
0	0	0	0	0	0	16	80.0
0	0	0	0	0	0	17	85.0
0	0	0	0	0	0	18	90.0
0	0	0	0	0	0	19	95.0
0	0	0	0	0	0	20	100.0

¹Intra-residue restraints, ²Sequential restraints, ³Medium range restraints, ⁴Long range restraints, ⁵Inter-chain restraints, ⁶ Number of models with violations

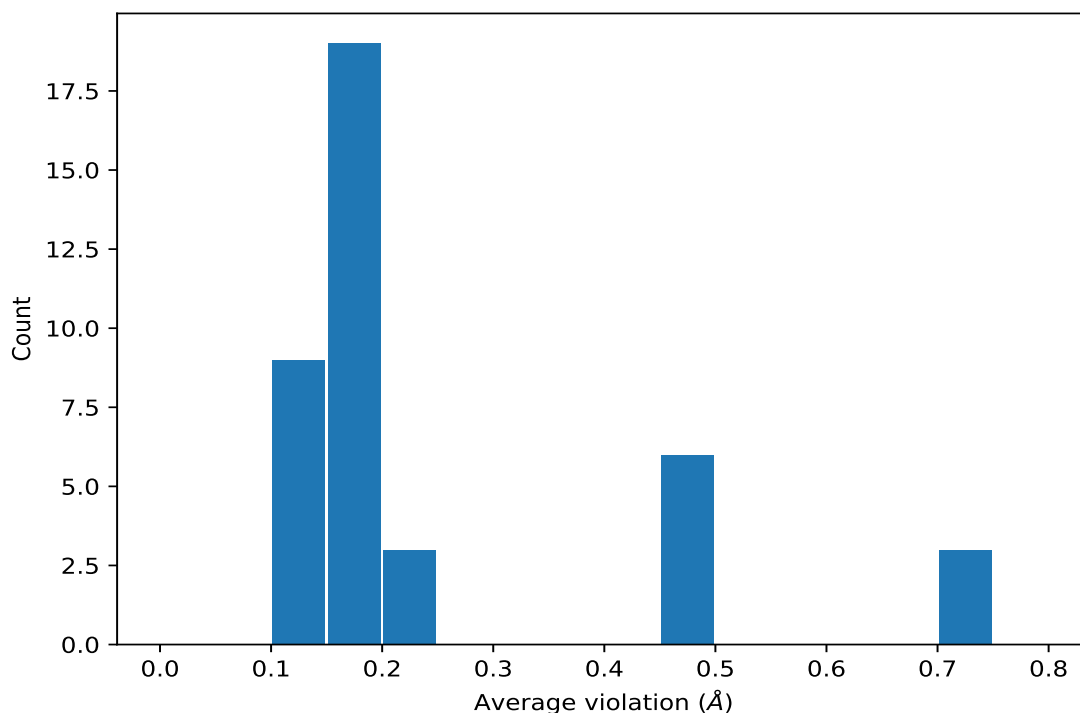
9.3.1 Bar graph : Distance violation statistics for the ensemble [i](#)



9.4 Most violated distance restraints in the ensemble [i](#)

9.4.1 Histogram : Distribution of mean distance violations [i](#)

The following histogram shows the distribution of the average value of the violation. The average is calculated for each restraint that is violated in more than one model over all the violated models in the ensemble



9.4.2 Table: Most violated distance restraints [i](#)

The following table provides the mean and the standard deviation of the violation for each restraint sorted by number of violated models and the mean value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	20	0.16	0.04	0.15
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	15	0.18	0.03	0.17
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	12	0.15	0.03	0.14
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	9	0.24	0.08	0.26
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	9	0.16	0.05	0.16
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	8	0.15	0.03	0.14
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	7	0.12	0.01	0.11
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	6	0.2	0.03	0.2
(1,21)	1:A:37:LEU:O	1:A:41:LEU:H	5	0.15	0.01	0.15
(1,1)	1:A:21:VAL:H	1:A:29:VAL:O	4	0.21	0.07	0.2
(1,25)	1:A:39:GLU:O	1:A:43:GLN:H	4	0.17	0.04	0.16
(1,20)	1:A:57:THR:O	1:A:67:TRP:N	4	0.16	0.05	0.14
(1,2)	1:A:21:VAL:N	1:A:29:VAL:O	3	0.19	0.04	0.16
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG21	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG22	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG23	3	0.16	0.01	0.16

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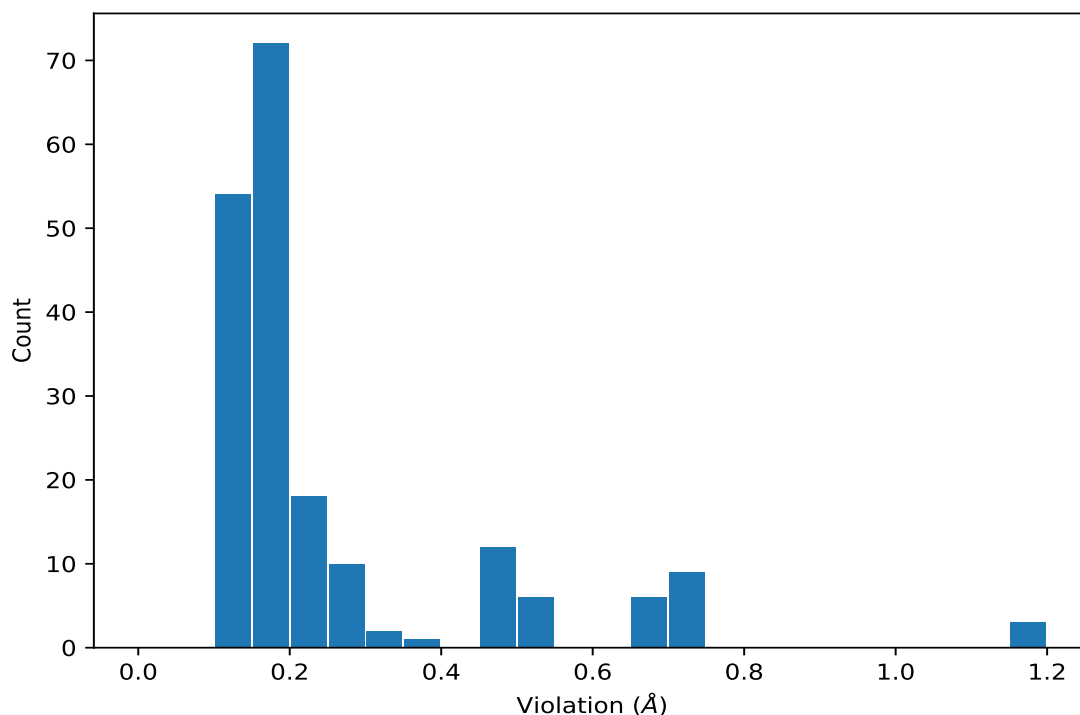
Key	Atom-1	Atom-2	Models ¹	Mean (Å)	SD ¹ (Å)	Median (Å)
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG21	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG22	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG23	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG21	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG22	3	0.16	0.01	0.16
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG23	3	0.16	0.01	0.16
(1,23)	1:A:38:GLU:O	1:A:42:ARG:H	3	0.13	0.01	0.13
(2,599)	1:A:29:VAL:HG11	1:B:163:VAL:HB	2	0.72	0.47	0.72
(2,599)	1:A:29:VAL:HG12	1:B:163:VAL:HB	2	0.72	0.47	0.72
(2,599)	1:A:29:VAL:HG13	1:B:163:VAL:HB	2	0.72	0.47	0.72
(2,752)	1:A:75:LEU:HD21	1:A:38:GLU:HB2	2	0.49	0.0	0.49
(2,752)	1:A:75:LEU:HD21	1:A:38:GLU:HB3	2	0.49	0.0	0.49
(2,752)	1:A:75:LEU:HD22	1:A:38:GLU:HB2	2	0.49	0.0	0.49
(2,752)	1:A:75:LEU:HD22	1:A:38:GLU:HB3	2	0.49	0.0	0.49
(2,752)	1:A:75:LEU:HD23	1:A:38:GLU:HB2	2	0.49	0.0	0.49
(2,752)	1:A:75:LEU:HD23	1:A:38:GLU:HB3	2	0.49	0.0	0.49
(2,795)	1:A:19:LEU:HD21	1:A:55:ILE:HA	2	0.18	0.04	0.18
(2,795)	1:A:19:LEU:HD22	1:A:55:ILE:HA	2	0.18	0.04	0.18
(2,795)	1:A:19:LEU:HD23	1:A:55:ILE:HA	2	0.18	0.04	0.18
(1,9)	1:A:18:PHE:O	1:A:56:ILE:H	2	0.16	0.01	0.16
(1,29)	1:A:41:LEU:O	1:A:45:ILE:H	2	0.15	0.02	0.15
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG11	2	0.13	0.02	0.13
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG12	2	0.13	0.02	0.13
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG13	2	0.13	0.02	0.13

¹Number of violated models, ²Standard deviation

9.5 All violated distance restraints [i](#)

9.5.1 Histogram : Distribution of distance violations [i](#)

The following histogram shows the distribution of the absolute value of the violation for all violated restraints in the ensemble.



9.5.2 Table : All distance violations [i](#)

The following table lists the absolute value of the violation for each restraint in the ensemble sorted by its value. The Key (restraint list ID, restraint ID) is the unique identifier for a given restraint. Rows with same key represent combinatorial or ambiguous restraints and are counted as a single restraint.

Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,599)	1:A:29:VAL:HG11	1:B:163:VAL:HB	6	1.19
(2,599)	1:A:29:VAL:HG12	1:B:163:VAL:HB	6	1.19
(2,599)	1:A:29:VAL:HG13	1:B:163:VAL:HB	6	1.19
(2,754)	1:A:78:MET:HE1	1:A:37:LEU:HD11	5	0.72
(2,754)	1:A:78:MET:HE1	1:A:37:LEU:HD12	5	0.72
(2,754)	1:A:78:MET:HE1	1:A:37:LEU:HD13	5	0.72
(2,754)	1:A:78:MET:HE2	1:A:37:LEU:HD11	5	0.72
(2,754)	1:A:78:MET:HE2	1:A:37:LEU:HD12	5	0.72
(2,754)	1:A:78:MET:HE2	1:A:37:LEU:HD13	5	0.72
(2,754)	1:A:78:MET:HE3	1:A:37:LEU:HD11	5	0.72
(2,754)	1:A:78:MET:HE3	1:A:37:LEU:HD12	5	0.72
(2,754)	1:A:78:MET:HE3	1:A:37:LEU:HD13	5	0.72
(2,561)	1:A:56:ILE:HD11	1:A:18:PHE:HD1	3	0.66
(2,561)	1:A:56:ILE:HD11	1:A:18:PHE:HD2	3	0.66
(2,561)	1:A:56:ILE:HD12	1:A:18:PHE:HD1	3	0.66
(2,561)	1:A:56:ILE:HD12	1:A:18:PHE:HD2	3	0.66

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,561)	1:A:56:ILE:HD13	1:A:18:PHE:HD1	3	0.66
(2,561)	1:A:56:ILE:HD13	1:A:18:PHE:HD2	3	0.66
(2,710)	1:A:21:VAL:HG11	1:A:31:LYS:HB2	2	0.53
(2,710)	1:A:21:VAL:HG11	1:A:31:LYS:HB3	2	0.53
(2,710)	1:A:21:VAL:HG12	1:A:31:LYS:HB2	2	0.53
(2,710)	1:A:21:VAL:HG12	1:A:31:LYS:HB3	2	0.53
(2,710)	1:A:21:VAL:HG13	1:A:31:LYS:HB2	2	0.53
(2,710)	1:A:21:VAL:HG13	1:A:31:LYS:HB3	2	0.53
(2,752)	1:A:75:LEU:HD21	1:A:38:GLU:HB2	11	0.49
(2,752)	1:A:75:LEU:HD21	1:A:38:GLU:HB3	11	0.49
(2,752)	1:A:75:LEU:HD22	1:A:38:GLU:HB2	11	0.49
(2,752)	1:A:75:LEU:HD22	1:A:38:GLU:HB3	11	0.49
(2,752)	1:A:75:LEU:HD23	1:A:38:GLU:HB2	11	0.49
(2,752)	1:A:75:LEU:HD23	1:A:38:GLU:HB3	11	0.49
(2,752)	1:A:75:LEU:HD21	1:A:38:GLU:HB2	16	0.49
(2,752)	1:A:75:LEU:HD21	1:A:38:GLU:HB3	16	0.49
(2,752)	1:A:75:LEU:HD22	1:A:38:GLU:HB2	16	0.49
(2,752)	1:A:75:LEU:HD22	1:A:38:GLU:HB3	16	0.49
(2,752)	1:A:75:LEU:HD23	1:A:38:GLU:HB2	16	0.49
(2,752)	1:A:75:LEU:HD23	1:A:38:GLU:HB3	16	0.49
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	16	0.35
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	1	0.31
(1,1)	1:A:21:VAL:H	1:A:29:VAL:O	17	0.31
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	18	0.28
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	7	0.27
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	14	0.26
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	6	0.26
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	12	0.26
(2,599)	1:A:29:VAL:HG11	1:B:163:VAL:HB	3	0.25
(2,599)	1:A:29:VAL:HG12	1:B:163:VAL:HB	3	0.25
(2,599)	1:A:29:VAL:HG13	1:B:163:VAL:HB	3	0.25
(1,2)	1:A:21:VAL:N	1:A:29:VAL:O	17	0.25
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	15	0.25
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	20	0.24
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	19	0.24
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	1	0.24
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	16	0.24
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	17	0.23
(1,25)	1:A:39:GLU:O	1:A:43:GLN:H	3	0.23
(1,20)	1:A:57:THR:O	1:A:67:TRP:N	6	0.23
(2,795)	1:A:19:LEU:HD21	1:A:55:ILE:HA	18	0.22
(2,795)	1:A:19:LEU:HD22	1:A:55:ILE:HA	18	0.22

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,795)	1:A:19:LEU:HD23	1:A:55:ILE:HA	18	0.22
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	16	0.22
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	8	0.22
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	8	0.22
(1,1)	1:A:21:VAL:H	1:A:29:VAL:O	8	0.22
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	7	0.21
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	14	0.21
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	7	0.2
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	18	0.2
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	17	0.19
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	20	0.19
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	13	0.19
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	19	0.19
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	1	0.19
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	10	0.19
(1,25)	1:A:39:GLU:O	1:A:43:GLN:H	9	0.19
(1,1)	1:A:21:VAL:H	1:A:29:VAL:O	18	0.19
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	10	0.18
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	5	0.18
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	12	0.18
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	5	0.18
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	12	0.18
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	12	0.18
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	14	0.18
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG21	18	0.17
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG22	18	0.17
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG23	18	0.17
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG21	18	0.17
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG22	18	0.17
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG23	18	0.17
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG21	18	0.17
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG22	18	0.17
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG23	18	0.17
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	7	0.17
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	16	0.17
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	9	0.17
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	11	0.17
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	12	0.17
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	15	0.17
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	8	0.17
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG21	4	0.16
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG22	4	0.16

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG23	4	0.16
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG21	4	0.16
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG22	4	0.16
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG23	4	0.16
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG21	4	0.16
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG22	4	0.16
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG23	4	0.16
(1,9)	1:A:18:PHE:O	1:A:56:ILE:H	17	0.16
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	14	0.16
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	18	0.16
(1,29)	1:A:41:LEU:O	1:A:45:ILE:H	12	0.16
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	2	0.16
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	6	0.16
(1,21)	1:A:37:LEU:O	1:A:41:LEU:H	1	0.16
(1,21)	1:A:37:LEU:O	1:A:41:LEU:H	18	0.16
(1,2)	1:A:21:VAL:N	1:A:29:VAL:O	8	0.16
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	19	0.16
(1,12)	1:A:53:ILE:O	1:A:71:ALA:N	15	0.16
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG21	2	0.15
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG22	2	0.15
(2,793)	1:A:20:VAL:HG11	1:A:52:THR:HG23	2	0.15
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG21	2	0.15
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG22	2	0.15
(2,793)	1:A:20:VAL:HG12	1:A:52:THR:HG23	2	0.15
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG21	2	0.15
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG22	2	0.15
(2,793)	1:A:20:VAL:HG13	1:A:52:THR:HG23	2	0.15
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG11	20	0.15
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG12	20	0.15
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG13	20	0.15
(1,9)	1:A:18:PHE:O	1:A:56:ILE:H	10	0.15
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	13	0.15
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	1	0.15
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	3	0.15
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	16	0.15
(1,26)	1:A:39:GLU:O	1:A:43:GLN:N	3	0.15
(1,21)	1:A:37:LEU:O	1:A:41:LEU:H	6	0.15
(1,21)	1:A:37:LEU:O	1:A:41:LEU:H	7	0.15
(1,2)	1:A:21:VAL:N	1:A:29:VAL:O	18	0.15
(2,795)	1:A:19:LEU:HD21	1:A:55:ILE:HA	10	0.14
(2,795)	1:A:19:LEU:HD22	1:A:55:ILE:HA	10	0.14
(2,795)	1:A:19:LEU:HD23	1:A:55:ILE:HA	10	0.14

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	14	0.14
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	15	0.14
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	20	0.14
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	1	0.14
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	4	0.14
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	13	0.14
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	15	0.14
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	18	0.14
(1,23)	1:A:38:GLU:O	1:A:42:ARG:H	17	0.14
(1,20)	1:A:57:THR:O	1:A:67:TRP:N	12	0.14
(1,20)	1:A:57:THR:O	1:A:67:TRP:N	19	0.14
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	10	0.13
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	12	0.13
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	18	0.13
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	6	0.13
(1,29)	1:A:41:LEU:O	1:A:45:ILE:H	1	0.13
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	10	0.13
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	12	0.13
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	7	0.13
(1,25)	1:A:39:GLU:O	1:A:43:GLN:H	20	0.13
(1,23)	1:A:38:GLU:O	1:A:42:ARG:H	12	0.13
(1,22)	1:A:37:LEU:O	1:A:41:LEU:N	1	0.13
(1,21)	1:A:37:LEU:O	1:A:41:LEU:H	12	0.13
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	11	0.12
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	9	0.12
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	4	0.12
(1,3)	1:A:21:VAL:O	1:A:29:VAL:H	6	0.12
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	5	0.12
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	9	0.12
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	17	0.12
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	20	0.12
(1,25)	1:A:39:GLU:O	1:A:43:GLN:H	4	0.12
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	20	0.12
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	10	0.12
(1,11)	1:A:53:ILE:O	1:A:71:ALA:H	20	0.12
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG11	3	0.11
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG12	3	0.11
(2,746)	1:A:54:ASP:HA	1:A:68:VAL:HG13	3	0.11
(1,4)	1:A:21:VAL:O	1:A:29:VAL:N	9	0.11
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	3	0.11
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	5	0.11
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	15	0.11

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Key	Atom-1	Atom-2	Model ID	Violation (Å)
(1,33)	1:A:43:GLN:O	1:A:47:GLN:H	18	0.11
(1,28)	1:A:40:PHE:O	1:A:44:LEU:N	2	0.11
(1,27)	1:A:40:PHE:O	1:A:44:LEU:H	11	0.11
(1,23)	1:A:38:GLU:O	1:A:42:ARG:H	11	0.11
(1,20)	1:A:57:THR:O	1:A:67:TRP:N	8	0.11
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	9	0.11
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	13	0.11
(1,19)	1:A:57:THR:O	1:A:67:TRP:H	17	0.11
(1,1)	1:A:21:VAL:H	1:A:29:VAL:O	10	0.11

10 Dihedral-angle violation analysis

Dihedral angle analysis failed due to data error in the dihedral angle restraints, possibly missing target value